

# Block diagonalization using similarity renormalization group flow equations

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By choosing appropriate generators for the Similarity Renormalization Group (SRG) flow equations, different patterns of decoupling in a Hamiltonian can be achieved. Sharp and smooth block-diagonal forms of phase-shift equivalent nucleon-nucleon potentials in momentum space are generated as examples and compared to analogous low-momentum interactions (“ $V_{\text{low } k}$ ”).

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The Similarity Renormalization Group (SRG) [1–3] applied to internucleon interactions is a continuous series of unitary transformations implemented as a flow equation for the evolving Hamiltonian  $H_s$ ,

$$\frac{dH_s}{ds} = [\eta_s, H_s] = [[G_s, H_s], H_s]. \quad (1)$$

Here  $s$  is a flow parameter and the flow operator  $G_s$  specifies the type of SRG [4]. Decoupling between low-energy and high-energy matrix elements is naturally achieved in a momentum basis by choosing a momentum-diagonal flow operator such as the kinetic energy  $T_{\text{rel}}$  or the diagonal of  $H_s$ ; either drives the Hamiltonian toward *band-diagonal* form. This decoupling leads to dramatically improved variational convergence in few-body nuclear systems compared to unevolved phenomenological or chiral effective field theory (EFT) potentials [5,6].

Renormalization Group (RG) methods that evolve  $NN$  interactions with a sharp or smooth cutoff in relative momentum, known generically as  $V_{\text{low } k}$ , rely on the invariance of the two-nucleon T matrix [7,8]. These approaches achieve a *block-diagonal* form characterized by a cutoff  $\Lambda$  (see left plots in Figs. 1 and 2). As implemented in Refs. [7,8], the high-momentum matrix elements are defined to be zero, but this is not required.

Block-diagonal decoupling of the sharp  $V_{\text{low } k}$  form can be generated using SRG flow equations by choosing a block-diagonal flow operator (see Refs. [9,10] for details),

$$G_s = \begin{pmatrix} PH_sP & 0 \\ 0 & QH_sQ \end{pmatrix} \equiv H_s^{\text{bd}}, \quad (2)$$

with projection operators  $P$  and  $Q = 1 - P$ . In a partial-wave momentum representation,  $P$  and  $Q$  are step functions defined by a sharp cutoff  $\Lambda$  on relative momenta. This choice for  $G_s$ , which means that  $\eta_s$  is nonzero only where  $G_s$  is zero, suppresses off-diagonal matrix elements such that the Hamiltonian approaches a block-diagonal form as  $s$  increases. If one considers a measure of the off-diagonal coupling of the Hamiltonian,

$$\text{Tr}[(QH_sP)^\dagger(QH_sP)] = \text{Tr}[PH_sQH_sP] \geq 0, \quad (3)$$

then its derivative is easily evaluated by applying the SRG equation, Eq. (1):

$$\begin{aligned} \frac{d}{ds} \text{Tr}[PH_sQH_sP] &= \text{Tr}[P\eta_sQ(QH_sQH_sP - QH_sPH_sP)] \\ &\quad + \text{Tr}[(PH_sPH_sQ - PH_sQH_sQ)Q\eta_sP] \\ &= -2 \text{Tr}[(Q\eta_sP)^\dagger(Q\eta_sP)] \leq 0. \end{aligned} \quad (4)$$

Thus, the off-diagonal  $QH_sP$  block will decrease in general as  $s$  increases [9,10].

The right plots in Figs. 1 and 2 result from evolving the  $N^3\text{LO}$  potential from Ref. [11] using the block-diagonal  $G_s$  of Eq. (2) with  $\Lambda = 2 \text{ fm}^{-1}$  until  $\lambda \equiv 1/s^{1/4} = 0.5 \text{ fm}^{-1}$ . The parameter  $\lambda$  is a quantitative measure of the progress toward block diagonalization made by the SRG evolution. The agreement between  $V_{\text{low } k}$  and SRG potentials for momenta below  $\Lambda$  is striking. A similar degree of universality is found in the other partial waves. Deriving an explicit connection between these approaches is the topic of an ongoing investigation.

The evolution with  $\lambda$  of two representative partial waves ( $^3S_1$  and  $^1P_1$ ) are shown in Figs. 3 and 4. The evolution of the “off-diagonal” matrix elements (meaning those outside the  $PH_sP$  and  $QH_sQ$  blocks) can be roughly understood from the dominance of the kinetic energy on the diagonal. Let the indices  $p$  and  $q$  run over indices of the momentum states in the  $P$  and  $Q$  spaces, respectively. To good approximation we can replace  $PH_sP$  and  $QH_sQ$  by their eigenvalues  $E_p$  and  $E_q$  in the SRG equations, yielding [9,10]

$$\frac{d}{ds} h_{pq} \approx \eta_{pq} E_q - E_p \eta_{pq} = -(E_p - E_q) \eta_{pq} \quad (5)$$

and

$$\eta_{pq} \approx E_p h_{pq} - h_{pq} E_q = (E_p - E_q) h_{pq}. \quad (6)$$

Combining these two results, we have the evolution of any off-diagonal matrix element:

$$\frac{d}{ds} h_{pq} \approx -(E_p - E_q)^2 h_{pq}. \quad (7)$$

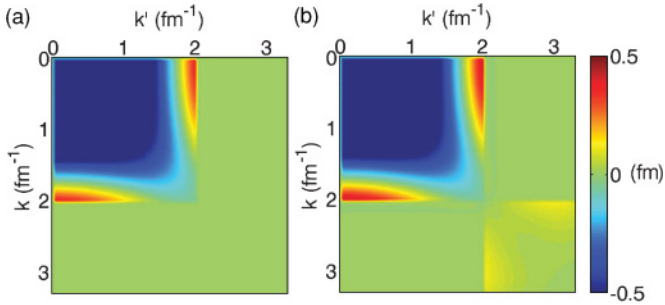


FIG. 1. (Color online) Comparison of momentum-space (a)  $V_{\text{low},k}$  and (b) SRG block-diagonal potentials with  $\Lambda = 2 \text{ fm}^{-1}$  evolved from an  $N^3\text{LO } ^3\text{S}_1$  potential [11].

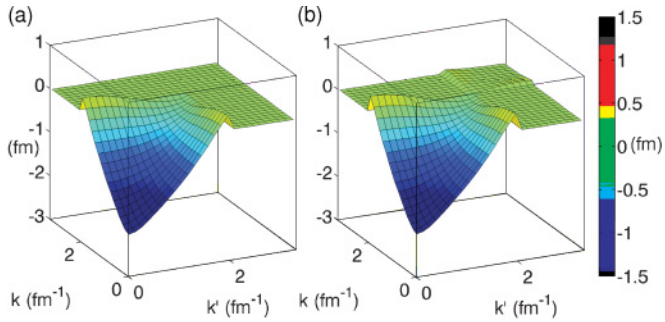


FIG. 2. (Color online) Comparison of momentum-space (a)  $V_{\text{low},k}$  and (b) SRG block-diagonal potentials with  $\Lambda = 2 \text{ fm}^{-1}$  evolved from an  $N^3\text{LO } ^3\text{S}_1$  potential [11].

In the  $NN$  case we can replace the eigenvalues by those for the relative kinetic energy, giving an explicit solution

$$h_{pq}(s) \approx h_{pq}(0)e^{-s(\epsilon_p - \epsilon_q)^2} \quad (8)$$

with  $\epsilon_p \equiv p^2/M$ . Thus the off-diagonal elements go to zero with the energy differences just like with the SRG with  $T_{\text{rel}}$ ; one can see the width of order  $1/\sqrt{s} = \lambda^2$  in the  $k^2$  plots of the evolving potential in Figs. 3 and 4.

While in principle the evolution to a sharp block-diagonal form means going to  $s = \infty (\lambda = 0)$ , in practice we need only take  $s$  as large as needed to quantitatively achieve the decoupling implied by Eq. (8). Furthermore, it should hold for more general definitions of  $P$  and  $Q$ . To smooth out the cutoff, we can introduce a smooth regulator  $f_\Lambda$ , which we take here to be an exponential form:

$$f_\Lambda(k) = e^{-(k^2/\Lambda^2)^n}, \quad (9)$$

with  $n$  an integer. For  $V_{\text{low},k}$  potentials, typical values used are  $n = 4$  and  $n = 8$  (the latter is considerably sharper but still numerically robust). By replacing  $H_s^{\text{bd}}$  with

$$G_s = f_\Lambda H_s f_\Lambda + (1 - f_\Lambda) H_s (1 - f_\Lambda), \quad (10)$$

we get a smooth block-diagonal potential.

A representative example with  $\Lambda = 2 \text{ fm}^{-1}$  and  $n = 4$  is shown in Fig. 5. We can evolve to  $\lambda = 1.5 \text{ fm}^{-1}$  without a problem. For smaller  $\lambda$  the overlap of the  $P$  and  $Q$  spaces becomes significant and the potential becomes distorted. This distortion indicates that there is no further benefit to evolving

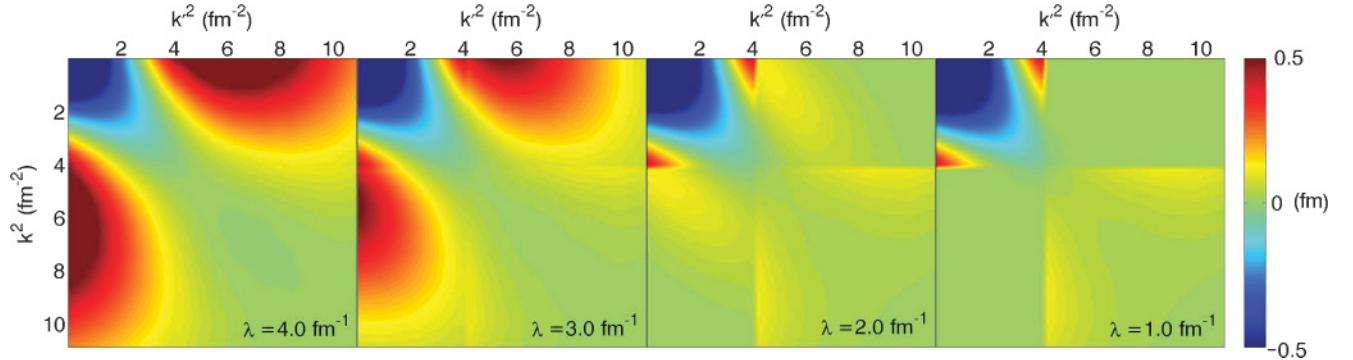


FIG. 3. (Color online) Evolution of the  $^3\text{S}_1$  partial wave with a sharp block-diagonal flow equation with  $\Lambda = 2 \text{ fm}^{-1}$  at  $\lambda = 4, 3, 2$ , and  $1 \text{ fm}^{-1}$ . The initial  $N^3\text{LO}$  potential is from Ref. [11].

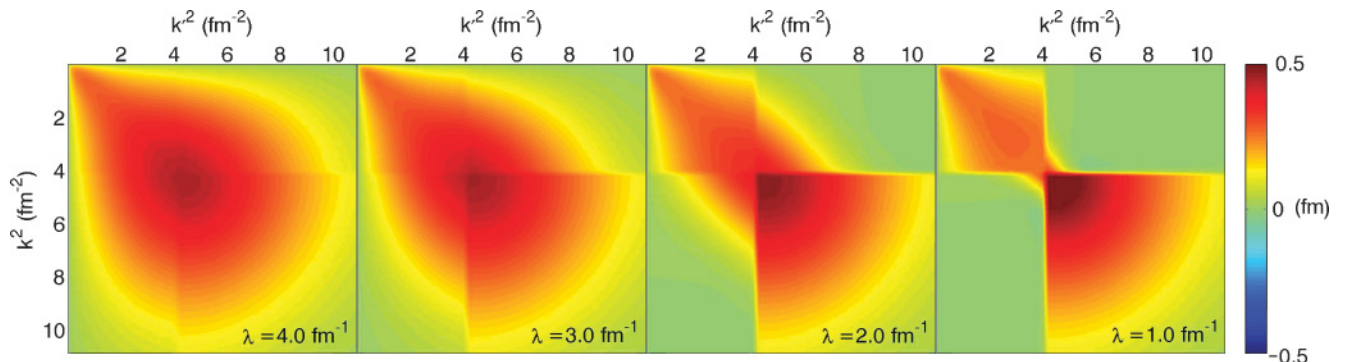


FIG. 4. (Color online) Same as Fig. 3 but for the  $^1\text{P}_1$  partial wave.

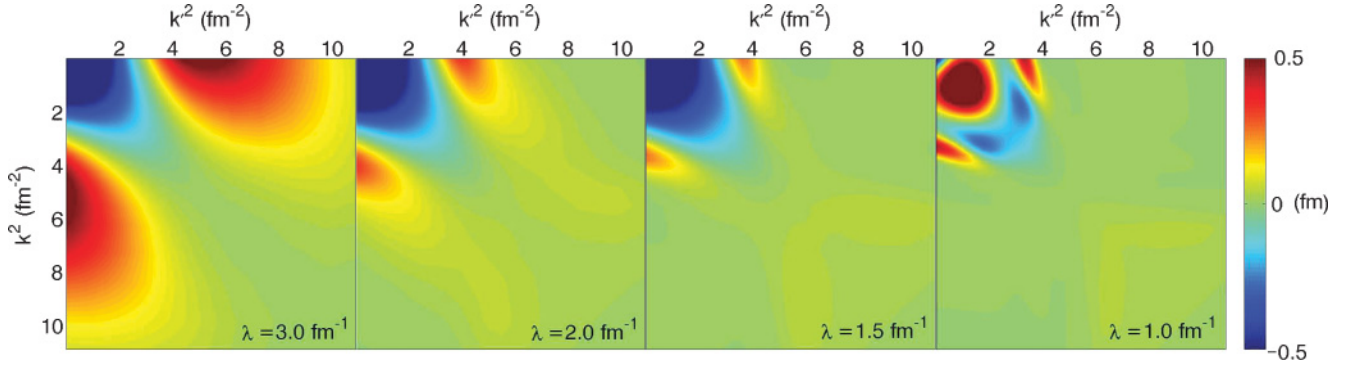


FIG. 5. (Color online) Evolution of the  $^3S_1$  partial wave with a smooth ( $n = 4$ ) block-diagonal flow equation with  $\Lambda = 2.0 \text{ fm}^{-1}$ , starting with the  $N^3\text{LO}$  potential from Ref. [11]. The flow parameter  $\lambda$  is 3, 2, 1.5, and  $1 \text{ fm}^{-1}$ .

in  $\lambda$  very far below  $\Lambda$ ; in fact the decoupling worsens for  $\lambda < \Lambda$  with a smooth regulator.

Another type of SRG that is second-order exact and yields similar block diagonalization is defined by

$$\eta_s = [T, P V_s Q + Q V_s P], \quad (11)$$

which can be implemented with  $P \rightarrow f_\Lambda$  and  $Q \rightarrow (1 - f_\Lambda)$ , with  $f_\Lambda$  either sharp or smooth. We can also consider bizarre choices for  $f_\Lambda$  in Eq. (10), such as defining it to be zero out to  $\Lambda_{\text{lower}}$ , then unity out to  $\Lambda$ , and then zero above that. This means that  $1 - f_\Lambda$  defines both low and high-momentum blocks and the region that is driven to zero consists of several rectangles. Results for two partial waves starting from the Argonne  $v_{18}$  potential [12] are shown in Fig. 6. Despite

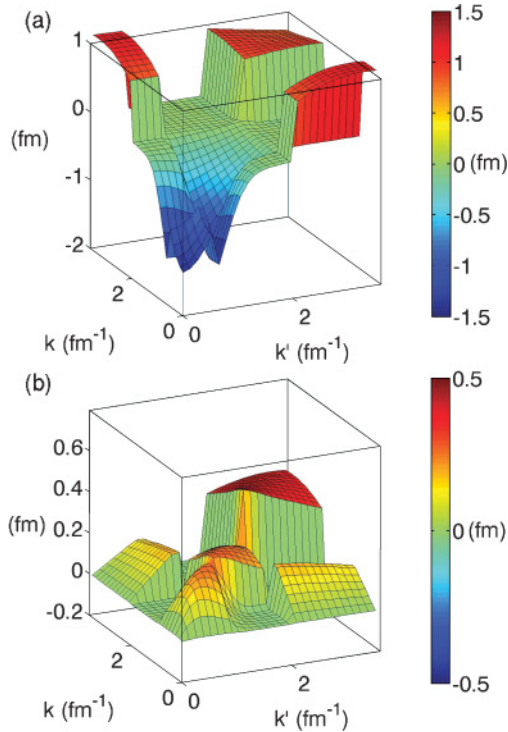


FIG. 6. (Color online) Evolved SRG potentials starting from Argonne  $v_{18}$  in the (a)  $^1S_0$  and (b)  $^1P_1$  partial waves to  $\lambda = 1 \text{ fm}^{-1}$  using a bizarre choice for  $G_s$  (see text).

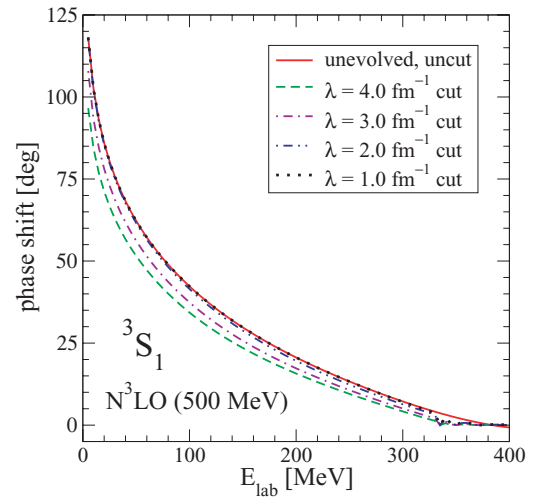


FIG. 7. (Color online) Phase shifts for the  $^3S_1$  partial wave from an initial  $N^3\text{LO}$  potential and the evolved sharp SRG block-diagonal potential with  $\Lambda = 2 \text{ fm}^{-1}$  at various  $\lambda$ , in each case with the potential set identically to zero above  $\Lambda$ .

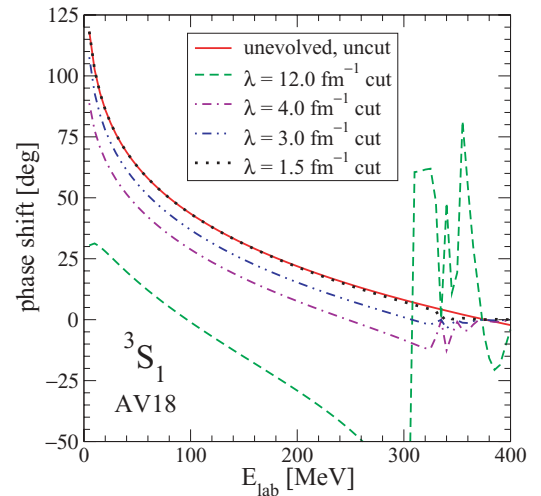


FIG. 8. (Color online) Same as Fig. 7 but with Argonne  $v_{18}$  as the initial potential [12].

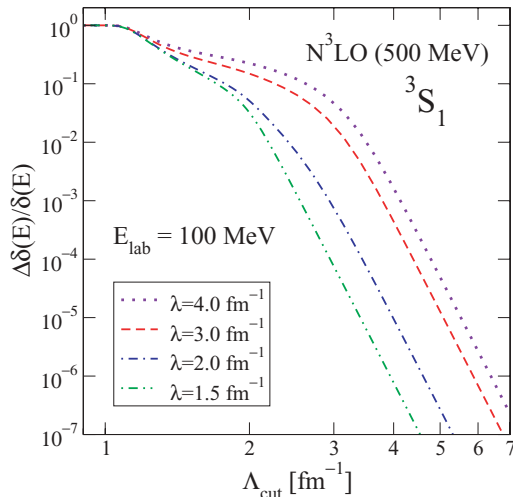


FIG. 9. (Color online) Errors in the phase shift at  $E_{\text{lab}} = 100$  MeV for the evolved sharp SRG block-diagonal potential with  $\Lambda = 2 \text{ fm}^{-1}$  for a range of  $\lambda$ 's and a regulator with  $n = 8$ .

the strange appearance, these remain unitary transformations of the original potential, with phase shifts and other  $NN$  observables the same as with the original potential. These choices provide a proof-of-principle that the decoupled regions can be tailored to the physics problem at hand.

Definitive tests of decoupling for  $NN$  observables are now possible for  $V_{\text{low } k}$  potentials since the unitary transformation of the SRG guarantees that no physics is lost. For example, in Figs. 7 and 8 we show  $^3S_1$  phase shifts from an SRG sharp block diagonalization with  $\Lambda = 2 \text{ fm}^{-1}$  for two different potentials. The phase shifts are calculated with the potentials cut sharply at  $\Lambda$ . That is, the matrix elements of the potential are set to zero above that point. The improved decoupling as

$\lambda$  decreases is evident in each case. By  $\lambda = 1 \text{ fm}^{-1}$  in Fig. 7, the unevolved and evolved curves are indistinguishable to the width of the line up to about 300 MeV.

In Fig. 9 we show a quantitative analysis of the decoupling as in Ref. [13]. The figure shows the relative error of the phase shift at 100 MeV calculated with a potential that is cut off by a smooth regulator as in Eq. (9) at a series of values  $\Lambda_{\text{cut}}$ . We observe the same universal decoupling behavior seen in Ref. [13]: a shoulder indicating the perturbative decoupling region, where the slope matches the power  $2n$  fixed by the smooth regulator. The onset of the shoulder in  $\Lambda_{\text{cut}}$  decreases with  $\lambda$  until it saturates for  $\lambda$  somewhat below  $\Lambda$ , leaving the shoulder at  $\Lambda_{\text{cut}} \approx \Lambda$ . Thus, as  $\lambda \rightarrow 0$  the decoupling scale is set by the cutoff  $\Lambda$ .

In the more conventional SRG, where we use  $\eta_s = [T, H_s] = [T, V_s]$ , it is easy to see that the evolution of the two-body potential in the two-particle system can be carried over directly to the three-particle system. In particular, it follows that the three-body potential does not depend on disconnected two-body parts [4,14]. If we could implement  $\eta_s$  as proposed here with analogous properties, we would have a tractable method for generating  $V_{\text{low } k}$  three-body forces. While it seems possible to define Fock-space operators with projectors  $P$  and  $Q$  that will not have problems with disconnected parts, it is not yet clear whether full decoupling in the few-body space can be realized. Work on this problem is in progress.

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